

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110723\
 Data File : BN028493.D
 Acq On : 07 Nov 2023 16:21
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN110723

Quant Time: Nov 08 00:30:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 08 00:27:19 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.717	152	7925	0.400	ng	0.00
7) Naphthalene-d8	10.509	136	24388	0.400	ng	0.00
13) Acenaphthene-d10	14.364	164	15437	0.400	ng	0.00
19) Phenanthrene-d10	17.109	188	36520	0.400	ng	0.00
29) Chrysene-d12	21.329	240	31348	0.400	ng	0.00
35) Perylene-d12	23.643	264	32534	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.291	112	9422	0.398	ng	0.00
5) Phenol-d6	6.887	99	12005	0.394	ng	0.00
8) Nitrobenzene-d5	8.865	82	8442	0.388	ng	0.00
11) 2-Methylnaphthalene-d10	12.102	152	16567	0.392	ng	0.00
14) 2,4,6-Tribromophenol	15.856	330	3031	0.350	ng	0.00
15) 2-Fluorobiphenyl	12.985	172	26637	0.393	ng	0.00
27) Fluoranthene-d10	19.153	212	42438	0.380	ng	0.00
31) Terphenyl-d14	19.766	244	31734	0.381	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.196	88	3882	0.407	ng	98
3) n-Nitrosodimethylamine	3.514	42	4357	0.398	ng	98
6) bis(2-Chloroethyl)ether	7.147	93	10032	0.398	ng	100
9) Naphthalene	10.552	128	30355	0.399	ng	100
10) Hexachlorobutadiene	10.851	225	5411	0.400	ng	# 100
12) 2-Methylnaphthalene	12.178	142	20436	0.392	ng	98
16) Acenaphthylene	14.076	152	30850	0.380	ng	99
17) Acenaphthene	14.418	154	21042	0.389	ng	99
18) Fluorene	15.412	166	29812	0.390	ng	100
20) 4,6-Dinitro-2-methylph...	15.497	198	1945	0.406	ng	99
21) 4-Bromophenyl-phenylether	16.315	248	9263	0.380	ng	98
22) Hexachlorobenzene	16.414	284	9868	0.387	ng	99
23) Atrazine	16.588	200	8109	0.377	ng	99
24) Pentachlorophenol	16.762	266	3593	0.328	ng	99
25) Phenanthrene	17.147	178	46723	0.386	ng	100
26) Anthracene	17.246	178	41801	0.377	ng	100
28) Fluoranthene	19.185	202	55477	0.382	ng	100
30) Pyrene	19.548	202	57261	0.384	ng	100
32) Benzo(a)anthracene	21.311	228	49652	0.379	ng	100
33) Chrysene	21.365	228	50925	0.389	ng	99
34) Bis(2-ethylhexyl)phtha...	21.266	149	27210	0.336	ng	99
36) Indeno(1,2,3-cd)pyrene	26.017	276	51883	0.395	ng	99
37) Benzo(b)fluoranthene	22.942	252	51170	0.382	ng	100
38) Benzo(k)fluoranthene	22.991	252	51197	0.391	ng	99
39) Benzo(a)pyrene	23.541	252	44529	0.389	ng	99
40) Dibenzo(a,h)anthracene	26.044	278	40488	0.396	ng	95
41) Benzo(g,h,i)perylene	26.748	276	45494	0.401	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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