

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062723\
 Data File : BN026045.D
 Acq On : 27 Jun 2023 20:01
 Operator : MA/JU
 Sample : 03376-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TW-2

Quant Time: Jun 28 05:50:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN062723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 28 05:23:31 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	5484	0.400	ng	0.00
7) Naphthalene-d8	10.782	136	16137	0.400	ng	0.00
13) Acenaphthene-d10	14.609	164	10090	0.400	ng	0.00
19) Phenanthrene-d10	17.356	188	20730	0.400	ng	0.00
29) Chrysene-d12	21.542	240	11565	0.400	ng	0.00
35) Perylene-d12	24.016	264	10424	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.528	112	2026	0.152	ng	0.00
5) Phenol-d6	7.146	99	1492	0.088	ng	0.00
8) Nitrobenzene-d5	9.159	82	3818	0.277	ng	0.01
11) 2-Methylnaphthalene-d10	12.374	152	7427	0.315	ng	0.00
14) 2,4,6-Tribromophenol	16.102	330	1353	0.369	ng	0.00
15) 2-Fluorobiphenyl	13.240	172	15873	0.407	ng	0.00
27) Fluoranthene-d10	19.384	212	18516	0.422	ng	0.00
31) Terphenyl-d14	19.974	244	16805	0.568	ng	0.00
Target Compounds						
						Qvalue
9) Naphthalene	10.835	128	3165	0.072	ng	# 88
10) Hexachlorobutadiene	11.113	225	532	0.063	ng	# 99
12) 2-Methylnaphthalene	12.449	142	2376	0.077	ng	95
16) Acenaphthylene	14.331	152	1411	0.030	ng	# 96
17) Acenaphthene	14.673	154	1165	0.038	ng	97
18) Fluorene	15.656	166	2457	0.062	ng	100
21) 4-Bromophenyl-phenylether	16.549	248	696	0.059	ng	# 85
22) Hexachlorobenzene	16.661	284	638	0.054	ng	# 94
25) Phenanthrene	17.393	178	4731	0.078	ng	98
26) Anthracene	17.493	178	2771	0.051	ng	98
28) Fluoranthene	19.416	202	3386	0.053	ng	# 95
30) Pyrene	19.779	202	2661	0.042	ng	99
32) Benzo(a)anthracene	21.533	228	1384	0.033	ng	# 81
33) Chrysene	21.578	228	1550	0.033	ng	# 82
34) Bis(2-ethylhexyl)phtha...	21.435	149	6452	0.174	ng	98
36) Indeno(1,2,3-cd)pyrene	26.610	276	1442	0.034	ng	# 73

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

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