

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071224\
 Data File : BN032632.D
 Acq On : 11 Jul 2024 12:53
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV243

Quant Time: Jul 12 01:41:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.570	152	4855	0.400	ng/ul	0.00
4) Naphthalene-d8	10.354	136	15125	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.216	164	9758	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.982	188	20125	0.400	ng/ul	0.00
17) Chrysene-d12	21.197	240	14911	0.400	ng/ul	0.00
23) Perylene-d12	23.395	264	14997	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.106	96	2332	0.400	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.971	152	8812	0.385	ng/ul	0.00
18) Fluoranthene-d10	19.022	212	20134	0.414	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.139	88	2629	0.396	ng/ul	98
5) Naphthalene	10.404	128	16564	0.408	ng/ul	99
7) 2-Methylnaphthalene	12.048	142	10701	0.422	ng/ul	99
8) 1-Methylnaphthalene	12.252	142	9849	0.376	ng/ul	99
10) Acenaphthylene	13.939	152	16487	0.418	ng/ul	100
11) Acenaphthene	14.276	153	13154	0.403	ng/ul	99
12) Fluorene	15.280	166	15098	0.405	ng/ul	100
14) Pentachlorophenol	16.666	266	1181	0.334	ng/ul	98
15) Phenanthrene	17.020	178	23031	0.420	ng/ul	99
16) Anthracene	17.126	178	19934	0.445	ng/ul	99
19) Fluoranthene	19.050	202	25819	0.426	ng/ul	99
20) Pyrene	19.417	202	27311	0.435	ng/ul	99
21) Benzo(a)anthracene	21.185	228	17943	0.378	ng/ul	100
22) Chrysene	21.235	228	27645	0.403	ng/ul	100
24) Benzo(b)fluoranthene	22.734	252	21408	0.405	ng/ul	99
25) Benzo(k)fluoranthene	22.775	252	24372	0.378	ng/ul	99
26) Benzo(a)pyrene	23.304	252	20257	0.442	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.623	276	24594	0.397	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.636	278	19086	0.397	ng/ul	100
29) Benzo(g,h,i)perylene	26.300	276	20533	0.396	ng/ul	99

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

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