

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041223\
 Data File : BN024852.D
 Acq On : 12 Apr 2023 17:40
 Operator : CG/JU
 Sample : PB152056BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS056

Quant Time: Apr 13 02:48:32 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN040723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 13 02:44:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	6912	0.400	ng/u1	0.00
4) Naphthalene-d8	10.765	136	19227	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.593	164	11148	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.338	188	24350	0.400	ng/u1	0.00
17) Chrysene-d12	21.535	240	19920	0.400	ng/u1	0.00
23) Perylene-d12	23.993	264	17306	0.400	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	6022	0.798	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.354	152	11079	0.449	ng/u1	0.00
18) Fluoranthene-d10	19.368	212	23204	0.469	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	15773	1.897	ng/u1	95
5) Naphthalene	10.814	128	20512	0.419	ng/u1	100
7) 2-Methylnaphthalene	12.425	142	13336	0.438	ng/u1	100
8) 1-Methylnaphthalene	12.640	142	13996	0.433	ng/u1	99
10) Acenaphthylene	14.316	152	17066	0.408	ng/u1	99
11) Acenaphthene	14.653	153	14892	0.416	ng/u1	100
12) Fluorene	15.643	166	16595	0.413	ng/u1	97
14) Pentachlorophenol	16.992	266	4457	0.858	ng/u1	99
15) Phenanthrene	17.381	178	27448	0.400	ng/u1	99
16) Anthracene	17.473	178	22424	0.393	ng/u1	99
19) Fluoranthene	19.400	202	30959	0.451	ng/u1	99
20) Pyrene	19.763	202	31589	0.449	ng/u1	98
21) Benzo(a)anthracene	21.520	228	27723	0.441	ng/u1	99
22) Chrysene	21.573	228	30846	0.453	ng/u1	100
24) Benzo(b)fluoranthene	23.242	252	31679	0.456	ng/u1	99
25) Benzo(k)fluoranthene	23.288	252	30463	0.450	ng/u1	98
26) Benzo(a)pyrene	23.888	252	25381	0.441	ng/u1	93
27) Indeno(1,2,3-cd)pyrene	26.561	276	28682	0.422	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.581	278	22055	0.413	ng/u1	97
29) Benzo(g,h,i)perylene	27.349	276	23820	0.415	ng/u1	99

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

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