

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023481.D
 Acq On : 26 Dec 2022 11:52
 Operator : CG/JU
 Sample : PB149682BS
 Misc :
 ALS Vial : 118 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS682

Quant Time: Dec 27 03:48:43 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Dec 25 05:17:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	6870	0.400	ng/u1	0.00
4) Naphthalene-d8	10.799	136	22546	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.631	164	12543	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.382	188	26195	0.400	ng/u1	0.00
17) Chrysene-d12	21.569	240	14558	0.400	ng/u1	0.00
23) Perylene-d12	24.013	264	11246	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	3820	0.512	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.388	152	11222	0.322	ng/u1	0.00
18) Fluoranthene-d10	19.411	212	21144	0.402	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.374	88	10356	1.356	ng/u1	90
5) Naphthalene	10.848	128	20913	0.316	ng/u1	100
7) 2-Methylnaphthalene	12.460	142	13197	0.312	ng/u1	100
8) 1-Methylnaphthalene	12.680	142	14260	0.327	ng/u1	100
10) Acenaphthylene	14.354	152	17803	0.320	ng/u1	100
11) Acenaphthene	14.696	153	14709	0.314	ng/u1	98
12) Fluorene	15.682	166	16229	0.308	ng/u1	99
14) Pentachlorophenol	17.031	266	3502	0.482	ng/u1	99
15) Phenanthrene	17.420	178	26495	0.311	ng/u1	99
16) Anthracene	17.513	178	24283	0.343	ng/u1	100
19) Fluoranthene	19.439	202	27584	0.391	ng/u1	99
20) Pyrene	19.801	202	28432	0.399	ng/u1	99
21) Benzo(a)anthracene	21.552	228	19880	0.351	ng/u1	99
22) Chrysene	21.607	228	20181	0.350	ng/u1	99
24) Benzo(b)fluoranthene	23.264	252	18438	0.317	ng/u1	98
25) Benzo(k)fluoranthene	23.311	252	17341	0.318	ng/u1	99
26) Benzo(a)pyrene	23.902	252	15255	0.331	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.567	276	18854	0.351	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.587	278	14882	0.357	ng/u1	99
29) Benzo(g,h,i)perylene	27.348	276	15929	0.364	ng/u1	98

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

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