

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110624\
 Data File : BM048479.D
 Acq On : 06 Nov 2024 16:13
 Operator : RC/JU
 Sample : PB164673BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS673

Quant Time: Nov 06 16:49:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 14:39:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	3191	0.400	ng/ul	0.00
4) Naphthalene-d8	10.475	136	9955	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.331	164	5397	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	9898	0.400	ng/ul	0.00
17) Chrysene-d12	21.269	240	8713	0.400	ng/ul	0.00
23) Perylene-d12	23.501	264	9585	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	3107	0.807	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.080	152	5378	0.418	ng/ul	0.00
18) Fluoranthene-d10	19.114	212	10796	0.441	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.168	88	8077	1.825	ng/ul#	67
5) Naphthalene	10.524	128	9335	0.374	ng/ul	98
7) 2-Methylnaphthalene	12.157	142	5731	0.365	ng/ul	96
8) 1-Methylnaphthalene	12.366	142	5825	0.362	ng/ul	100
10) Acenaphthylene	14.053	152	7743	0.378	ng/ul	100
11) Acenaphthene	14.391	153	6209	0.379	ng/ul	99
12) Fluorene	15.395	166	6704	0.376	ng/ul	98
14) Pentachlorophenol	16.753	266	2798	0.770	ng/ul	99
15) Phenanthrene	17.129	178	10034	0.386	ng/ul	99
16) Anthracene	17.230	178	9047	0.374	ng/ul	99
19) Fluoranthene	19.146	202	12744	0.395	ng/ul	98
20) Pyrene	19.509	202	13878	0.396	ng/ul	98
21) Benzo(a)anthracene	21.260	228	10504	0.368	ng/ul	99
22) Chrysene	21.306	228	15482	0.452	ng/ul	100
24) Benzo(b)fluoranthene	22.832	252	12549	0.420	ng/ul	98
25) Benzo(k)fluoranthene	22.876	252	16009	0.457	ng/ul	99
26) Benzo(a)pyrene	23.414	252	12471	0.428	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.769	276	17432	0.428	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.786	278	13531	0.427	ng/ul	99
29) Benzo(g,h,i)perylene	26.450	276	15141	0.443	ng/ul	98

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

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