

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122724\
 Data File : BM049291.D
 Acq On : 28 Dec 2024 13:56
 Operator : RC/JU
 Sample : PB165854BS
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS854

Quant Time: Dec 29 21:57:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM121824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 26 11:35:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.544	152	5748	0.400	ng/ul	0.00
4) Naphthalene-d8	10.303	136	16715	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.182	164	9341	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.933	188	17942	0.400	ng/ul	0.00
17) Chrysene-d12	21.135	240	17469	0.400	ng/ul	0.00
23) Perylene-d12	23.288	264	20870	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	3699	0.558	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.903	152	7125	0.317	ng/ul	0.00
18) Fluoranthene-d10	18.968	212	14722	0.297	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.168	88	10544	1.361	ng/ul#	78
5) Naphthalene	10.352	128	12324	0.295	ng/ul	98
7) 2-Methylnaphthalene	11.980	142	7590	0.286	ng/ul	100
8) 1-Methylnaphthalene	12.200	142	7727	0.285	ng/ul	100
10) Acenaphthylene	13.899	152	11116	0.287	ng/ul	98
11) Acenaphthene	14.242	153	8091	0.286	ng/ul	99
12) Fluorene	15.241	166	8843	0.277	ng/ul	98
14) Pentachlorophenol	16.595	266	3083	0.622	ng/ul	96
15) Phenanthrene	16.975	178	13425	0.286	ng/ul	99
16) Anthracene	17.064	178	12851	0.299	ng/ul	98
19) Fluoranthene	18.996	202	17144	0.271	ng/ul	100
20) Pyrene	19.363	202	18577	0.276	ng/ul	98
21) Benzo(a)anthracene	21.120	228	17224	0.283	ng/ul	99
22) Chrysene	21.170	228	19293	0.296	ng/ul	97
24) Benzo(b)fluoranthene	22.648	252	18697	0.280	ng/ul	92
25) Benzo(k)fluoranthene	22.692	252	20235	0.277	ng/ul#	90
26) Benzo(a)pyrene	23.195	252	18508	0.289	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	25.424	276	26350	0.291	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.441	278	20403	0.290	ng/ul#	94
29) Benzo(g,h,i)perylene	26.068	276	21681	0.298	ng/ul	96

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

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