

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032962.D
 Acq On : 10 Nov 2021 20:58
 Operator : CG/JU
 Sample : PB140517BS
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS517

Quant Time: Nov 11 08:23:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 13:31:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.469	152	1497	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	5762	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.117	164	3279	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.862	188	6772	0.400	ng/ul	0.00
17) Chrysene-d12	21.049	240	5527	0.400	ng/ul	0.00
23) Perylene-d12	23.159	264	4428	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.864	96	1537	0.804	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.836	152	2906	0.355	ng/ul	0.00
18) Fluoranthene-d10	18.892	212	6540	0.390	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.898	88	4225	2.152	ng/ul#	90
5) Naphthalene	10.284	128	6273	0.375	ng/ul	100
7) 2-Methylnaphthalene	11.912	142	3915	0.338	ng/ul	100
8) 1-Methylnaphthalene	12.133	142	3782	0.333	ng/ul	100
10) Acenaphthylene	13.837	152	4892	0.343	ng/ul	98
11) Acenaphthene	14.182	153	4578	0.368	ng/ul	99
12) Fluorene	15.175	166	5185	0.353	ng/ul	97
14) Pentachlorophenol	16.539	266	886	0.502	ng/ul	98
15) Phenanthrene	16.903	178	8238	0.366	ng/ul	99
16) Anthracene	16.994	178	6677	0.338	ng/ul	99
19) Fluoranthene	18.923	202	9695	0.374	ng/ul	98
20) Pyrene	19.285	202	10135	0.376	ng/ul	99
21) Benzo(a)anthracene	21.035	228	7085	0.347	ng/ul	100
22) Chrysene	21.088	228	8826	0.384	ng/ul	99
24) Benzo(b)fluoranthene	22.536	252	8089	0.398	ng/ul	96
25) Benzo(k)fluoranthene	22.575	252	8783	0.403	ng/ul	94
26) Benzo(a)pyrene	23.067	252	6489	0.388	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	25.220	276	7768	0.403	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.223	278	6191	0.406	ng/ul	96
29) Benzo(g,h,i)perylene	25.850	276	7051	0.422	ng/ul	99

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

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