

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102721\
 Data File : BM032764.D
 Acq On : 27 Oct 2021 17:13
 Operator : CG/JU
 Sample : PB140207BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS207

Quant Time: Oct 27 17:45:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM102721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 27 15:49:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.539	152	2673	0.400	ng/ul	0.00
4) Naphthalene-d8	10.316	136	10276	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.190	164	6083	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.925	188	13006	0.400	ng/ul	0.00
17) Chrysene-d12	21.101	240	12397	0.400	ng/ul	0.00
23) Perylene-d12	23.227	264	11419	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.910	96	2384	0.684	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.913	152	4955	0.348	ng/ul	0.00
18) Fluoranthene-d10	18.950	212	12235	0.360	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.944	88	6305	1.713	ng/ul#	88
5) Naphthalene	10.366	128	10316	0.336	ng/ul	99
7) 2-Methylnaphthalene	11.990	142	6604	0.322	ng/ul	100
8) 1-Methylnaphthalene	12.210	142	6294	0.313	ng/ul	100
10) Acenaphthylene	13.910	152	7478	0.322	ng/ul	98
11) Acenaphthene	14.251	153	7640	0.325	ng/ul	99
12) Fluorene	15.241	166	8700	0.321	ng/ul	98
14) Pentachlorophenol	16.599	266	2340	0.539	ng/ul	99
15) Phenanthrene	16.963	178	13978	0.322	ng/ul	99
16) Anthracene	17.057	178	11248	0.314	ng/ul	100
19) Fluoranthene	18.981	202	17076	0.332	ng/ul	100
20) Pyrene	19.343	202	17597	0.331	ng/ul	98
21) Benzo(a)anthracene	21.087	228	14399	0.328	ng/ul	99
22) Chrysene	21.137	228	17906	0.342	ng/ul	99
24) Benzo(b)fluoranthene	22.597	252	17749	0.345	ng/ul	99
25) Benzo(k)fluoranthene	22.639	252	19379	0.344	ng/ul	99
26) Benzo(a)pyrene	23.138	252	14417	0.333	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.321	276	19060	0.333	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.323	278	15260	0.333	ng/ul	99
29) Benzo(g,h,i)perylene	25.965	276	17483	0.336	ng/ul	98

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

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