

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120423\
 Data File : BM043179.D
 Acq On : 04 Dec 2023 16:57
 Operator : MA/JU
 Sample : PB157549BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS549

Quant Time: Dec 04 17:35:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM112423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 02 23:54:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.847	152	1035	0.400	ng/ul	0.02
4) Naphthalene-d8	10.627	136	3423	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.450	164	1992	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.224	188	3897	0.400	ng/ul	0.00
17) Chrysene-d12	21.409	240	2386	0.400	ng/ul	0.00
23) Perylene-d12	23.753	264	3086	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.155	96	826	0.520	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.238	152	1399	0.305	ng/ul	0.02
18) Fluoranthene-d10	19.252	212	2730	0.314	ng/ul	0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.189	88	2291	1.343	ng/ul#	86
5) Naphthalene	10.677	128	2489	0.268	ng/ul	99
7) 2-Methylnaphthalene	12.304	142	1472	0.268	ng/ul	98
8) 1-Methylnaphthalene	12.508	142	1587	0.259	ng/ul	97
10) Acenaphthylene	14.186	152	2734	0.269	ng/ul	98
11) Acenaphthene	14.515	153	2111	0.274	ng/ul	98
12) Fluorene	15.528	166	2133	0.270	ng/ul	98
14) Pentachlorophenol	16.887	266	658	0.624	ng/ul	94
15) Phenanthrene	17.266	178	3091	0.278	ng/ul	99
16) Anthracene	17.380	178	3227	0.306	ng/ul	98
19) Fluoranthene	19.280	202	3951	0.292	ng/ul	99
20) Pyrene	19.647	202	4066	0.283	ng/ul	97
21) Benzo(a)anthracene	21.400	228	2438	0.315	ng/ul	99
22) Chrysene	21.447	228	4083	0.283	ng/ul	98
24) Benzo(b)fluoranthene	23.034	252	3201	0.307	ng/ul	99
25) Benzo(k)fluoranthene	23.084	252	4354	0.304	ng/ul	99
26) Benzo(a)pyrene	23.660	252	3368	0.296	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.222	276	4475	0.291	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.239	278	3386	0.285	ng/ul	95
29) Benzo(g,h,i)perylene	26.973	276	4097	0.274	ng/ul	99

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

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