

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042623\
 Data File : BM039696.D
 Acq On : 27 Apr 2023 01:44
 Operator : CG/JU
 Sample : PB152427BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS427

Quant Time: Apr 27 02:23:42 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM040623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 27 01:14:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.876	152	2117	0.400	ng/ul	0.02
4) Naphthalene-d8	10.645	136	11281	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.483	164	6506	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.255	188	9742	0.400	ng/ul	0.00
17) Chrysene-d12	21.438	240	7638	0.400	ng/ul	# 0.00
23) Perylene-d12	23.808	264	6258	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	3457	0.882	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	5258	0.403	ng/ul	0.02
18) Fluoranthene-d10	19.276	212	9370	0.394	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.244	88	8490	2.023	ng/ul#	77
5) Naphthalene	10.694	128	10035	0.374	ng/ul	97
7) 2-Methylnaphthalene	12.328	142	5580	0.370	ng/ul	97
8) 1-Methylnaphthalene	12.526	142	6154	0.360	ng/ul	99
10) Acenaphthylene	14.210	152	9669	0.383	ng/ul	98
11) Acenaphthene	14.544	153	7648	0.385	ng/ul	99
12) Fluorene	15.557	166	7803	0.394	ng/ul	98
14) Pentachlorophenol	16.951	266	1090	0.469	ng/ul#	83
15) Phenanthrene	17.297	178	10462	0.387	ng/ul	98
16) Anthracene	17.403	178	7784	0.378	ng/ul	100
19) Fluoranthene	19.309	202	11822	0.374	ng/ul	98
20) Pyrene	19.666	202	13506	0.373	ng/ul	99
21) Benzo(a)anthracene	21.429	228	7019	0.398	ng/ul	100
22) Chrysene	21.476	228	9521	0.441	ng/ul	99
24) Benzo(b)fluoranthene	23.083	252	8451	0.426	ng/ul	96
25) Benzo(k)fluoranthene	23.136	252	8439	0.446	ng/ul	96
26) Benzo(a)pyrene	23.709	252	8136	0.407	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.282	276	11180	0.444	ng/ul	98
28) Dibenzo(a,h)anthracene	26.302	278	8458	0.479	ng/ul	98
29) Benzo(g,h,i)perylene	27.019	276	10347	0.428	ng/ul	98

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

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