

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092721\
 Data File : BM032250.D
 Acq On : 28 Sep 2021 21:40
 Operator : CG/JU
 Sample : PB139277BS
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS277

Quant Time: Sep 29 03:25:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM092721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 27 13:37:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	6327	0.400	ng/ul	0.00
4) Naphthalene-d8	10.450	136	19030	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.303	164	8324	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.028	188	15634	0.400	ng/ul	0.00
17) Chrysene-d12	21.190	240	12541	0.400	ng/ul	0.00
23) Perylene-d12	23.355	264	10937	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.989	96	5409	0.795	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.038	152	9813	0.387	ng/ul	-0.01
18) Fluoranthene-d10	19.048	212	14782	0.434	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.024	88	14156	2.023	ng/ul	90
5) Naphthalene	10.495	128	21409	0.379	ng/ul	99
7) 2-Methylnaphthalene	12.115	142	13000	0.360	ng/ul	100
8) 1-Methylnaphthalene	12.335	142	12210	0.346	ng/ul	100
10) Acenaphthylene	14.019	152	14083	0.422	ng/ul	100
11) Acenaphthene	14.364	153	11801	0.384	ng/ul	99
12) Fluorene	15.346	166	12503	0.359	ng/ul	99
14) Pentachlorophenol	16.695	266	2836	0.717	ng/ul	99
15) Phenanthrene	17.066	178	18925	0.372	ng/ul	99
16) Anthracene	17.157	178	15682	0.374	ng/ul	99
19) Fluoranthene	19.075	202	20416	0.390	ng/ul	98
20) Pyrene	19.437	202	20491	0.390	ng/ul	100
21) Benzo(a)anthracene	21.175	228	17610	0.397	ng/ul	100
22) Chrysene	21.226	228	20270	0.391	ng/ul	100
24) Benzo(b)fluoranthene	22.708	252	20843	0.367	ng/ul	95
25) Benzo(k)fluoranthene	22.749	252	20314	0.370	ng/ul	95
26) Benzo(a)pyrene	23.258	252	16472	0.386	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.494	276	21586	0.396	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.499	278	17649	0.398	ng/ul	98
29) Benzo(g,h,i)perylene	26.155	276	20441	0.396	ng/ul	99

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

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