

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080521\
 Data File : BM031559.D
 Acq On : 07 Aug 2021 20:21
 Operator : CG/JU
 Sample : PB138086BS
 Misc :
 ALS Vial : 85 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS086

Quant Time: Aug 08 08:41:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 08 08:38:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	1261	0.400	ng/ul	0.00
4) Naphthalene-d8	10.347	136	4525	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.216	164	2408	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	4487	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.159	240	3479	0.400	ng/ul	# 0.00
23) Perylene-d12	23.319	264	3439	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	854	0.610	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.944	152	2274	0.298	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	3877	0.334	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.220	88	2162	1.528	ng/ul#	77
5) Naphthalene	10.397	128	4034	0.299	ng/ul	98
7) 2-Methylnaphthalene	12.020	142	2637	0.286	ng/ul	99
8) 1-Methylnaphthalene	12.241	142	2397	0.263	ng/ul	100
10) Acenaphthylene	13.941	152	3242	0.316	ng/ul	99
11) Acenaphthene	14.281	153	2559	0.296	ng/ul	100
12) Fluorene	15.278	166	2796	0.281	ng/ul	98
14) Pentachlorophenol	16.628	266	810	0.564	ng/ul	99
15) Phenanthrene	17.013	178	4257	0.299	ng/ul	99
16) Anthracene	17.103	178	4023	0.302	ng/ul	98
19) Fluoranthene	19.031	202	4762	0.321	ng/ul#	93
20) Pyrene	19.394	202	4852	0.322	ng/ul#	94
21) Benzo(a)anthracene	21.144	228	4455	0.317	ng/ul	97
22) Chrysene	21.195	228	4632	0.319	ng/ul	98
24) Benzo(b)fluoranthene	22.674	252	4761	0.307	ng/ul	91
25) Benzo(k)fluoranthene	22.716	252	4787	0.299	ng/ul#	92
26) Benzo(a)pyrene	23.224	252	4194	0.309	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.456	276	5354	0.305	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.468	278	4235	0.300	ng/ul#	95
29) Benzo(g,h,i)perylene	26.108	276	4604	0.308	ng/ul#	95

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

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