

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081121\
 Data File : BM031660.D
 Acq On : 11 Aug 2021 15:01
 Operator : CG/JU
 Sample : PB138247BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS247

Quant Time: Aug 11 15:36:48 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 : Wed Aug 11 11:28:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.576	152	1231	0.400	ng/ul	0.00
4) Naphthalene-d8	10.338	136	4715	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.212	164	2488	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.968	188	4456	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.161	240	3377	0.400	ng/ul	# 0.00
23) Perylene-d12	23.321	264	3359	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.186	96	738	0.540	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.935	152	2205	0.278	ng/ul	0.00
18) Fluoranthene-d10	18.997	212	3741	0.332	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.217	88	2015	1.459	ng/ul#	75
5) Naphthalene	10.388	128	4010	0.286	ng/ul	99
7) 2-Methylnaphthalene	12.011	142	2606	0.272	ng/ul	97
8) 1-Methylnaphthalene	12.232	142	2385	0.252	ng/ul	100
10) Acenaphthylene	13.932	152	3369	0.318	ng/ul	100
11) Acenaphthene	14.277	153	2570	0.288	ng/ul	99
12) Fluorene	15.270	166	2769	0.270	ng/ul	95
14) Pentachlorophenol	16.628	266	742	0.520	ng/ul	98
15) Phenanthrene	17.006	178	4159	0.294	ng/ul	97
16) Anthracene	17.100	178	3885	0.293	ng/ul	99
19) Fluoranthene	19.027	202	4583	0.318	ng/ul#	95
20) Pyrene	19.390	202	4714	0.322	ng/ul#	93
21) Benzo(a)anthracene	21.144	228	4276	0.314	ng/ul	97
22) Chrysene	21.198	228	4476	0.317	ng/ul	99
24) Benzo(b)fluoranthene	22.677	252	4459	0.295	ng/ul	85
25) Benzo(k)fluoranthene	22.720	252	4649	0.297	ng/ul#	88
26) Benzo(a)pyrene	23.227	252	4081	0.308	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	25.466	276	5118	0.299	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.478	278	4054	0.294	ng/ul#	90
29) Benzo(g,h,i)perylene	26.110	276	4450	0.304	ng/ul#	92

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

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