

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082622\
 Data File : BM036443.D
 Acq On : 27 Aug 2022 02:23
 Operator : CG/JU
 Sample : PB147244BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS244

Quant Time: Aug 27 03:26:28 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM082622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 14:26:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	3365	0.400	ng/ul	0.00
4) Naphthalene-d8	10.600	136	14013	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.436	164	7437	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.187	188	15641	0.400	ng/ul	0.00
17) Chrysene-d12	21.363	240	16278	0.400	ng/ul	0.00
23) Perylene-d12	23.692	264	12875	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	3723	0.822	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.206	152	7786	0.348	ng/ul	0.02
18) Fluoranthene-d10	19.210	212	16878	0.839	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.239	88	9466	2.212	ng/ul#	79
5) Naphthalene	10.650	128	13718	0.347	ng/ul	98
7) 2-Methylnaphthalene	12.277	142	8292	0.350	ng/ul	96
8) 1-Methylnaphthalene	12.486	142	8494	0.354	ng/ul	99
10) Acenaphthylene	14.163	152	11757	0.382	ng/ul	100
11) Acenaphthene	14.501	153	9391	0.378	ng/ul	99
12) Fluorene	15.496	166	10540	0.362	ng/ul	99
14) Pentachlorophenol	16.853	266	2005	0.649	ng/ul	98
15) Phenanthrene	17.225	178	17457	0.313	ng/ul	99
16) Anthracene	17.326	178	15393	0.347	ng/ul	99
19) Fluoranthene	19.243	202	20624	0.362	ng/ul	99
20) Pyrene	19.605	202	21578	0.365	ng/ul	97
21) Benzo(a)anthracene	21.351	228	17685	0.262	ng/ul	99
22) Chrysene	21.401	228	22668	0.245	ng/ul	98
24) Benzo(b)fluoranthene	22.979	252	25926	0.277	ng/ul	91
25) Benzo(k)fluoranthene	23.025	252	23977	0.254	ng/ul#	90
26) Benzo(a)pyrene	23.590	252	19037	0.321	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.105	276	22021	0.491	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.128	278	17518	0.461	ng/ul	91
29) Benzo(g,h,i)perylene	26.836	276	20098	0.496	ng/ul	97

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

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