

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040824\
 Data File : BM045067.D
 Acq On : 09 Apr 2024 07:52
 Operator : MA/JU
 Sample : PB160054BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS054

Quant Time: Apr 09 09:04:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM040424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 09 02:13:29 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.648	152	1504	0.400	ng/ul	0.00
4) Naphthalene-d8	10.430	136	4915	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.288	164	2624	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.056	188	5519	0.400	ng/ul	0.00
17) Chrysene-d12	21.260	240	4125	0.400	ng/ul	0.00
23) Perylene-d12	23.479	264	4857	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	1271	0.643	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.041	152	2420	0.363	ng/ul	0.00
18) Fluoranthene-d10	19.089	212	5010	0.481	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.243	88	3140	1.658	ng/ul#	80
5) Naphthalene	10.479	128	4459	0.342	ng/ul	97
7) 2-Methylnaphthalene	12.118	142	2647	0.340	ng/ul	99
8) 1-Methylnaphthalene	12.321	142	2794	0.329	ng/ul	97
10) Acenaphthylene	14.011	152	4593	0.351	ng/ul	97
11) Acenaphthene	14.353	153	3348	0.357	ng/ul	99
12) Fluorene	15.361	166	3642	0.347	ng/ul	99
14) Pentachlorophenol	16.722	266	925	0.715	ng/ul	96
15) Phenanthrene	17.098	178	5537	0.356	ng/ul	99
16) Anthracene	17.204	178	5511	0.360	ng/ul	95
19) Fluoranthene	19.117	202	6608	0.448	ng/ul	97
20) Pyrene	19.484	202	6914	0.444	ng/ul	97
21) Benzo(a)anthracene	21.252	228	5259	0.366	ng/ul	97
22) Chrysene	21.295	228	7282	0.406	ng/ul	99
24) Benzo(b)fluoranthene	22.812	252	6716	0.383	ng/ul	96
25) Benzo(k)fluoranthene	22.859	252	8187	0.417	ng/ul	93
26) Benzo(a)pyrene	23.385	252	6677	0.399	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.733	276	8669	0.378	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.756	278	6713	0.369	ng/ul#	88
29) Benzo(g,h,i)perylene	26.397	276	7825	0.395	ng/ul	97

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

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