

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070324\
 Data File : BN032488.D
 Acq On : 04 Jul 2024 23:34
 Operator : MA/JU
 Sample : PB161807BS
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS807

Quant Time: Jul 05 01:10:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	5439	0.400	ng/ul	0.00
4) Naphthalene-d8	10.376	136	15794	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.248	164	8470	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.008	188	15402	0.400	ng/ul	0.00
17) Chrysene-d12	21.211	240	14953	0.400	ng/ul	# 0.00
23) Perylene-d12	23.427	264	18331	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	4387	0.726	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.977	152	7814	0.330	ng/ul	0.00
18) Fluoranthene-d10	19.045	212	15048	0.334	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	11730	1.704	ng/ul	91
5) Naphthalene	10.426	128	14122	0.338	ng/ul	99
7) 2-Methylnaphthalene	12.048	142	8541	0.308	ng/ul	100
8) 1-Methylnaphthalene	12.268	142	8699	0.303	ng/ul	99
10) Acenaphthylene	13.966	152	12666	0.360	ng/ul	99
11) Acenaphthene	14.309	153	9400	0.344	ng/ul	99
12) Fluorene	15.308	166	9675	0.297	ng/ul	99
14) Pentachlorophenol	16.678	266	2684	0.770	ng/ul	100
15) Phenanthrene	17.050	178	14220	0.337	ng/ul	99
16) Anthracene	17.147	178	13176	0.350	ng/ul	97
19) Fluoranthene	19.073	202	17697	0.311	ng/ul#	95
20) Pyrene	19.440	202	19205	0.323	ng/ul	99
21) Benzo(a)anthracene	21.194	228	19113	0.354	ng/ul	96
22) Chrysene	21.246	228	20359	0.344	ng/ul	97
24) Benzo(b)fluoranthene	22.760	252	21253	0.267	ng/ul	85
25) Benzo(k)fluoranthene	22.804	252	22036	0.285	ng/ul#	86
26) Benzo(a)pyrene	23.333	252	20392	0.368	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.666	276	27341	0.315	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.673	278	21216	0.329	ng/ul#	89
29) Benzo(g,h,i)perylene	26.354	276	21447	0.310	ng/ul	92

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

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