

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072524\
 Data File : BN032984.D
 Acq On : 25 Jul 2024 22:34
 Operator : MA/JU
 Sample : PB162260BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS260

Quant Time: Jul 26 01:19:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 15:43:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.533	152	2554	0.400	ng/ul	0.00
4) Naphthalene-d8	10.306	136	9856	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.171	164	7075	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	16781	0.400	ng/ul	0.00
17) Chrysene-d12	21.154	240	16104	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	19390	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.069	96	2446	0.757	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.917	152	5680	0.353	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	17626	0.387	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.098	88	6927	1.881	ng/ul#	70
5) Naphthalene	10.361	128	10946	0.432	ng/ul	98
7) 2-Methylnaphthalene	11.994	142	6407	0.374	ng/ul	100
8) 1-Methylnaphthalene	12.198	142	6625	0.357	ng/ul	99
10) Acenaphthylene	13.898	152	10960	0.379	ng/ul	99
11) Acenaphthene	14.236	153	8833	0.406	ng/ul	99
12) Fluorene	15.244	166	9788	0.386	ng/ul	99
14) Pentachlorophenol	16.629	266	2082	0.609	ng/ul	99
15) Phenanthrene	16.983	178	15729	0.349	ng/ul	99
16) Anthracene	17.084	178	14544	0.359	ng/ul	99
19) Fluoranthene	19.014	202	20821	0.393	ng/ul	99
20) Pyrene	19.381	202	22516	0.396	ng/ul	99
21) Benzo(a)anthracene	21.139	228	20466	0.373	ng/ul	99
22) Chrysene	21.192	228	24904	0.388	ng/ul	100
24) Benzo(b)fluoranthene	22.685	252	22167	0.391	ng/ul	97
25) Benzo(k)fluoranthene	22.732	252	27519	0.405	ng/ul	95
26) Benzo(a)pyrene	23.255	252	23230	0.436	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.547	276	28755	0.355	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.550	278	22113	0.344	ng/ul	99
29) Benzo(g,h,i)perylene	26.228	276	22355	0.347	ng/ul	98

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

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