

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071624\
 Data File : BN032771.D
 Acq On : 17 Jul 2024 09:59
 Operator : MA/JU
 Sample : PB161927BS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS927

Quant Time: Jul 17 11:29:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 17 02:35:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	5155	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	13210	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.197	164	8403	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	15982	0.400	ng/ul	0.00
17) Chrysene-d12	21.173	240	15984	0.400	ng/ul	# 0.00
23) Perylene-d12	23.374	264	18776	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	4572	0.739	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.922	152	8816	0.442	ng/ul	-0.01
18) Fluoranthene-d10	19.003	212	15580	0.299	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.135	88	12327	1.748	ng/ul	90
5) Naphthalene	10.365	128	13077	0.369	ng/ul	99
7) 2-Methylnaphthalene	11.993	142	9608	0.434	ng/ul	98
8) 1-Methylnaphthalene	12.213	142	9675	0.423	ng/ul	100
10) Acenaphthylene	13.920	152	14966	0.441	ng/ul	100
11) Acenaphthene	14.262	153	10356	0.369	ng/ul	99
12) Fluorene	15.257	166	10678	0.333	ng/ul	99
14) Pentachlorophenol	16.636	266	2908	1.037	ng/ul	100
15) Phenanthrene	17.003	178	16286	0.374	ng/ul	98
16) Anthracene	17.096	178	15947	0.448	ng/ul	98
19) Fluoranthene	19.031	202	18349	0.282	ng/ul	96
20) Pyrene	19.398	202	20191	0.300	ng/ul	97
21) Benzo(a)anthracene	21.159	228	22346	0.439	ng/ul#	90
22) Chrysene	21.211	228	22799	0.310	ng/ul#	90
24) Benzo(b)fluoranthene	22.713	252	24200	0.365	ng/ul#	72
25) Benzo(k)fluoranthene	22.754	252	23816	0.295	ng/ul#	71
26) Benzo(a)pyrene	23.278	252	23290	0.406	ng/ul#	68
27) Indeno(1,2,3-cd)pyrene	25.592	276	30133	0.388	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.592	278	23591	0.392	ng/ul#	77
29) Benzo(g,h,i)perylene	26.273	276	23372	0.360	ng/ul#	80

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

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