

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027942.D
 Acq On : 29 Sep 2023 00:55
 Operator : MA/JU
 Sample : PB155830BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS830

Quant Time: Sep 29 04:01:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 27 03:52:19 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	4238	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	16771	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	11172	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	24625	0.400	ng/ul	0.00
17) Chrysene-d12	21.536	240	21541	0.400	ng/ul	0.00
23) Perylene-d12	24.006	264	19426	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	4949	0.927	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	14147	0.557	ng/ul	0.00
18) Fluoranthene-d10	19.376	212	35585	0.541	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.363	88	12821	2.308	ng/ul#	83
5) Naphthalene	10.828	128	20874	0.436	ng/ul	100
7) 2-Methylnaphthalene	12.434	142	14525	0.478	ng/ul	100
8) 1-Methylnaphthalene	12.649	142	14403	0.459	ng/ul	99
10) Acenaphthylene	14.328	152	24060	0.493	ng/ul	99
11) Acenaphthene	14.661	153	17214	0.408	ng/ul	98
12) Fluorene	15.647	166	19878	0.401	ng/ul	98
14) Pentachlorophenol	16.987	266	3861	0.670	ng/ul	99
15) Phenanthrene	17.396	178	32170	0.416	ng/ul	98
16) Anthracene	17.485	178	30734	0.474	ng/ul	99
19) Fluoranthene	19.408	202	38762	0.445	ng/ul	100
20) Pyrene	19.776	202	40675	0.453	ng/ul	98
21) Benzo(a)anthracene	21.519	228	37533	0.450	ng/ul	99
22) Chrysene	21.574	228	35958	0.418	ng/ul	99
24) Benzo(b)fluoranthene	23.240	252	33125	0.383	ng/ul	98
25) Benzo(k)fluoranthene	23.293	252	32999	0.393	ng/ul	98
26) Benzo(a)pyrene	23.895	252	29995	0.431	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.599	276	33893	0.380	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.623	278	27871	0.388	ng/ul	99
29) Benzo(g,h,i)perylene	27.414	276	27298	0.360	ng/ul	99

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

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