

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072023\
 Data File : BN026634.D
 Acq On : 21 Jul 2023 18:54
 Operator : MA/JU
 Sample : PB154137BS
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS137

Quant Time: Jul 22 00:52:10 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 21 19:02:16 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	1454	0.400	ng/ul	0.00
4) Naphthalene-d8	10.641	136	6001	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.472	164	3896	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.228	188	7494	0.400	ng/ul	0.00
17) Chrysene-d12	21.419	240	5168	0.400	ng/ul	0.00
23) Perylene-d12	23.790	264	5720	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.203	96	1590	0.830	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.242	152	3272	0.401	ng/ul	0.00
18) Fluoranthene-d10	19.255	212	8518	0.475	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.237	88	3443	1.900	ng/ul#	89
5) Naphthalene	10.691	128	5588	0.339	ng/ul	96
7) 2-Methylnaphthalene	12.313	142	3205	0.343	ng/ul	97
8) 1-Methylnaphthalene	12.522	142	4034	0.403	ng/ul	99
10) Acenaphthylene	14.199	152	6647	0.348	ng/ul	99
11) Acenaphthene	14.536	153	5152	0.354	ng/ul	98
12) Fluorene	15.531	166	5508	0.361	ng/ul	99
14) Pentachlorophenol	16.890	266	976	0.516	ng/ul	97
15) Phenanthrene	17.266	178	8390	0.352	ng/ul	99
16) Anthracene	17.363	178	7237	0.349	ng/ul	98
19) Fluoranthene	19.283	202	10164	0.420	ng/ul#	96
20) Pyrene	19.645	202	11190	0.425	ng/ul#	95
21) Benzo(a)anthracene	21.402	228	6789	0.357	ng/ul	100
22) Chrysene	21.452	228	8500	0.393	ng/ul	99
24) Benzo(b)fluoranthene	23.065	252	7792	0.350	ng/ul	96
25) Benzo(k)fluoranthene	23.112	252	8536	0.365	ng/ul#	97
26) Benzo(a)pyrene	23.688	252	7392	0.373	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.244	276	9745	0.375	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.258	278	7682	0.385	ng/ul	95
29) Benzo(g,h,i)perylene	27.002	276	8428	0.367	ng/ul	98

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

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