

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
 Data File : BP014772.D
 Acq On : 21 Apr 2023 00:05
 Operator : CG/JU
 Sample : 02340-16
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHE17

Quant Time: Apr 21 02:54:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 20 02:32:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.152	152	229830	20.000	ng/ul	0.00
20) Naphthalene-d8	10.975	136	920161	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.781	164	550243	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.527	188	1165334	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	1078732	20.000	ng/ul	0.00
88) Perylene-d12	24.168	264	1231239	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.487	96	14328	2.478	ng/uL	0.00
4) Pyridine-d5	3.922	84	93410	5.933	ng/ul	0.00
7) Phenol-d5	7.287	99	118237	5.950	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.463	67	362336	29.556	ng/ul	0.00
11) 2-Chlorophenol-d4	7.669	132	353435	23.949	ng/ul	0.00
15) 4-Methylphenol-d8	8.846	113	217979	14.071	ng/ul	0.00
21) Nitrobenzene-d5	9.322	128	200359	34.749	ng/ul	0.00
24) 2-Nitrophenol-d4	10.046	143	187079	35.923	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.581	165	379682	26.851	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.104	131	512479	23.745	ng/ul	0.00
46) Dimethylphthalate-d6	14.181	166	1397522	32.076	ng/ul	0.00
49) Acenaphthylene-d8	14.475	160	1573814	31.170	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	30649	4.423	ng/ul	0.00
60) Fluorene-d10	15.769	176	1225323	32.776	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	132720	28.143	ng/ul	0.00
73) Anthracene-d10	17.627	188	2185093	38.814	ng/ul	0.00
81) Pyrene-d10	19.839	212	2532004	40.974	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.004	264	2574284	40.340	ng/ul	-0.01

Target Compounds Qvalue

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

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