

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101723\
 Data File : BP017705.D
 Acq On : 17 Oct 2023 19:15
 Operator : MA/JU
 Sample : PB156344TB 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB344

Quant Time: Oct 18 00:45:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 16:14:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	89145	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	362728	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	229709	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.392	188	532633	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	554961	20.000	ng/u1	-0.01
88) Perylene-d12	24.098	264	641180	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	22371	9.252	ng/uL	-0.01
4) Pyridine-d5	3.693	84	290550	43.470	ng/u1	0.00
7) Phenol-d5	7.058	99	367109	44.233	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	240180	47.720	ng/u1	0.00
11) 2-Chlorophenol-d4	7.417	132	305008	47.478	ng/u1	0.00
15) 4-Methylphenol-d8	8.622	113	291238	43.934	ng/u1	0.00
21) Nitrobenzene-d5	9.111	128	144770	47.517	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	158256	46.189	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	276177	42.888	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	390507	42.308	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	939047	47.201	ng/u1	0.00
49) Acenaphthylene-d8	14.299	160	1003248	46.829	ng/u1	0.00
54) 4-Nitrophenol-d4	14.851	143	117534	35.325	ng/u1	0.00
60) Fluorene-d10	15.604	176	825068	48.754	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	112894	29.987	ng/u1	0.00
73) Anthracene-d10	17.492	188	1348103	49.028	ng/u1	0.00
81) Pyrene-d10	19.751	212	1677125	50.211	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.927	264	1828881	50.809	ng/u1	-0.01

Target Compounds	Qvalue

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

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