

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017794.D
 Acq On : 25 Oct 2023 01:04
 Operator : MA/JU
 Sample : PB156501BL
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK501

Quant Time: Oct 25 04:39:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	152749	20.000	ng/u1	0.00
20) Naphthalene-d8	10.710	136	618503	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.575	164	373026	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.369	188	778181	20.000	ng/u1	0.00
79) Chrysene-d12	21.510	240	630626	20.000	ng/u1	# 0.00
88) Perylene-d12	24.063	264	650910	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	24643	6.002	ng/uL	0.00
4) Pyridine-d5	3.687	84	302561	28.145	ng/u1	0.00
7) Phenol-d5	7.040	99	397184	31.247	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	263626	33.583	ng/u1	0.00
11) 2-Chlorophenol-d4	7.399	132	326337	31.478	ng/u1	-0.01
15) 4-Methylphenol-d8	8.605	113	316310	31.688	ng/u1	0.00
21) Nitrobenzene-d5	9.087	128	157312	31.388	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.804	143	175731	31.735	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	297405	27.379	ng/u1	0.00
31) 4-Chloroaniline-d4	10.887	131	385944	27.149	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	942478	30.097	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	1059699	30.345	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	108742	24.371	ng/u1	0.00
60) Fluorene-d10	15.581	176	793285	29.306	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.728	200	108997	21.264	ng/u1	0.00
73) Anthracene-d10	17.469	188	1200275	30.183	ng/u1	0.00
81) Pyrene-d10	19.727	212	1327598	33.579	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.892	264	1095102	30.207	ng/u1	0.00

Target Compounds Qvalue

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

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