

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017537.D
 Acq On : 04 Oct 2023 17:41
 Operator : MA/JU
 Sample : PB156006BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK006

Quant Time: Oct 05 00:46:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	70503	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	276816	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	170478	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.422	188	387869	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	383845	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	442762	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	16313	9.250	ng/uL	0.00
4) Pyridine-d5	3.728	84	228412	49.483	ng/u1	0.00
7) Phenol-d5	7.099	99	270984	45.683	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	173529	47.941	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	220079	47.119	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	212412	44.730	ng/u1	0.00
21) Nitrobenzene-d5	9.151	128	106722	50.113	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	117214	48.800	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	198533	44.584	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	290320	45.767	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	701369	52.743	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	731857	49.911	ng/u1	0.00
54) 4-Nitrophenol-d4	14.875	143	108781	46.713	ng/u1	0.00
60) Fluorene-d10	15.639	176	579211	51.352	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	103153	42.659	ng/u1	0.00
73) Anthracene-d10	17.522	188	942421	52.491	ng/u1	0.00
81) Pyrene-d10	19.774	212	1166935	54.711	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.962	264	1226237	54.879	ng/u1	-0.01

Target Compounds Qvalue

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

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