

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
 Data File : BP017724.D
 Acq On : 19 Oct 2023 11:17
 Operator : MA/JU
 Sample : PB156373BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK373

Quant Time: Oct 20 00:52:34 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	7.893	152	103066	20.000	ng/u1	0.00
20) Naphthalene-d8	10.728	136	396274	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.593	164	241940	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.387	188	518441	20.000	ng/u1	0.01
79) Chrysene-d12	21.527	240	431505	20.000	ng/u1	0.01
88) Perylene-d12	24.086	264	459968	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	20824	7.516	ng/uL	0.00
4) Pyridine-d5	3.699	84	251302	34.645	ng/u1	0.00
7) Phenol-d5	7.052	99	290892	33.917	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	199279	37.623	ng/u1	0.00
11) 2-Chlorophenol-d4	7.416	132	245386	35.079	ng/u1	0.00
15) 4-Methylphenol-d8	8.616	113	229963	34.143	ng/u1	0.00
21) Nitrobenzene-d5	9.105	128	113666	35.398	ng/u1	0.00
24) 2-Nitrophenol-d4	9.822	143	123993	34.948	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.346	165	215925	31.025	ng/u1	0.00
31) 4-Chloroaniline-d4	10.905	131	289684	31.806	ng/u1	0.01
46) Dimethylphthalate-d6	14.010	166	730528	35.969	ng/u1	0.00
49) Acenaphthylene-d8	14.293	160	780991	34.481	ng/u1	0.01
54) 4-Nitrophenol-d4	14.840	143	75596	26.122	ng/u1	0.01
60) Fluorene-d10	15.598	176	619097	35.263	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	92030	26.949	ng/u1	0.01
73) Anthracene-d10	17.487	188	938386	35.419	ng/u1	0.01
81) Pyrene-d10	19.745	212	1077219	39.819	ng/u1	0.01
92) Benzo(a)pyrene-d12	23.916	264	923883	36.063	ng/u1	0.02

Target Compounds Qvalue

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

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