

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017752.D
 Acq On : 23 Oct 2023 16:01
 Operator : MA/JU
 Sample : PB156470BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK470

Quant Time: Oct 24 05:14:07 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	148725	20.000	ng/u1	0.00
20) Naphthalene-d8	10.722	136	575328	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	370125	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	766952	20.000	ng/u1	0.00
79) Chrysene-d12	21.521	240	730004	20.000	ng/u1	0.00
88) Perylene-d12	24.080	264	819091	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	22930	5.736	ng/uL	0.00
4) Pyridine-d5	3.693	84	277851	26.545	ng/u1	0.00
7) Phenol-d5	7.052	99	324172	26.193	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	213955	27.993	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	281314	27.869	ng/u1	0.00
15) 4-Methylphenol-d8	8.610	113	253913	26.126	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	125236	26.863	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	141190	27.410	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	255189	25.256	ng/u1	0.00
31) 4-Chloroaniline-d4	10.898	131	321983	24.350	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	858515	27.631	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	923739	26.659	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	79411	17.937	ng/u1	0.00
60) Fluorene-d10	15.592	176	749879	27.920	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	114418	22.649	ng/u1	0.00
73) Anthracene-d10	17.480	188	1123603	28.668	ng/u1	0.00
81) Pyrene-d10	19.739	212	1336809	29.209	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.910	264	1315454	28.834	ng/u1	0.01

Target Compounds Qvalue

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

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