

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
 Data File : BP012595.D
 Acq On : 10 Nov 2022 12:40
 Operator : CG/JU
 Sample : PB148940BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK940

Quant Time: Nov 10 21:34:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 09 23:28:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	100453	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	470116	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.416	164	318610	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	709206	20.000	ng/u1	0.00
79) Chrysene-d12	21.280	240	748250	20.000	ng/u1	0.00
88) Perylene-d12	23.651	264	706630	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	32591	13.224	ng/uL	0.00
4) Pyridine-d5	3.646	84	190594	27.365	ng/u1	0.00
7) Phenol-d5	6.952	99	552769	65.276	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	352219	69.516	ng/u1	0.00
11) 2-Chlorophenol-d4	7.299	132	451020	69.211	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	445601	66.521	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	229725	66.041	ng/u1	0.00
24) 2-Nitrophenol-d4	9.652	143	253245	66.285	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	424910	61.734	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	239422	24.695	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	1520780	69.249	ng/u1	0.00
49) Acenaphthylene-d8	14.110	160	1669141	67.185	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	218683	63.243	ng/u1	0.00
60) Fluorene-d10	15.416	176	1299817	69.804	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	225749	57.551	ng/u1	0.00
73) Anthracene-d10	17.281	188	2195063	71.272	ng/u1	0.00
81) Pyrene-d10	19.522	212	2593657	61.856	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.498	264	2594104	74.470	ng/u1	0.00

Target Compounds Qvalue

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

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