

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012729.D
 Acq On : 25 Nov 2022 16:34
 Operator : CG/JU
 Sample : PB149032BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK032

Quant Time: Nov 26 06:21:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 25 15:59:22 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.028	152	442384	20.000 ng/u1	0.00
20) Naphthalene-d8	10.846	136	1950142	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.663	164	1351961	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.422	188	2953556	20.000 ng/u1	0.00
79) Chrysene-d12	21.504	240	2828242	20.000 ng/u1	0.00
88) Perylene-d12	24.027	264	2651380	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.399	96	63007	6.103 ng/uL	0.00
4) Pyridine-d5	3.817	84	862693	27.924 ng/u1	0.00
7) Phenol-d5	7.169	99	1076769	29.621 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	702711	31.371 ng/u1	0.00
11) 2-Chlorophenol-d4	7.552	132	870618	31.431 ng/u1	0.00
15) 4-Methylphenol-d8	8.722	113	887489	30.301 ng/u1	0.00
21) Nitrobenzene-d5	9.193	128	415763	34.833 ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	463410	36.565 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	900860	30.464 ng/u1	0.00
31) 4-Chloroaniline-d4	10.975	131	1297043	32.030 ng/u1	0.00
46) Dimethylphthalate-d6	14.069	166	3247958	34.903 ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	3513104	32.355 ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	462985	27.906 ng/u1	0.00
60) Fluorene-d10	15.657	176	2755732	33.814 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	373568	26.463 ng/u1	0.00
73) Anthracene-d10	17.516	188	4525992	34.812 ng/u1	0.00
81) Pyrene-d10	19.745	212	5159739	31.397 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.863	264	4561260	35.024 ng/u1	0.00

Target Compounds Qvalue

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

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