

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012217.D
 Acq On : 20 Oct 2022 17:06
 Operator : CG/JU
 Sample : PB148363BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK363

Quant Time: Oct 20 23:24:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	350271	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1461364	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.475	164	875421	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1693800	20.000	ng/u1	0.00
79) Chrysene-d12	21.327	240	1393950	20.000	ng/u1	0.00
88) Perylene-d12	23.733	264	1463213	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	59836	6.913	ng/uL	0.00
4) Pyridine-d5	3.681	84	485610	19.471	ng/u1	0.00
7) Phenol-d5	6.993	99	977901	32.339	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	605684	33.027	ng/u1	0.00
11) 2-Chlorophenol-d4	7.357	132	780943	33.928	ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	753958	32.080	ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	370463	34.248	ng/u1	0.00
24) 2-Nitrophenol-d4	9.716	143	400163	34.288	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	705889	32.245	ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	654638	21.047	ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	2110482	34.657	ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2409812	34.669	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	302799	26.542	ng/u1	0.00
60) Fluorene-d10	15.475	176	1833429	34.997	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	265968	27.602	ng/u1	0.00
73) Anthracene-d10	17.333	188	2695690	36.308	ng/u1	0.00
81) Pyrene-d10	19.574	212	2812231	35.434	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.580	264	2629501	35.835	ng/u1	0.00

Target Compounds Qvalue

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

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