

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010854.D
 Acq On : 26 Jun 2022 06:49
 Operator : CG/JU
 Sample : PB145767BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK767

Quant Time: Jun 26 22:57:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	414420	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.510	136	1646022	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.369	164	866967	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.133	188	1635118	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1634986	20.000	ng/u1	#-0.01
88) Perylene-d12	23.609	264	1739993	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	78527	7.168	ng/uL	0.00
4) Pyridine-d5	3.610	84	992953	34.049	ng/u1	0.00
7) Phenol-d5	6.893	99	1159873	35.356	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.057	67	741120	34.543	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	988606	36.322	ng/u1	0.00
15) 4-Methylphenol-d8	8.428	113	921865	35.548	ng/u1	0.00
21) Nitrobenzene-d5	8.875	128	441800	42.162	ng/u1	0.00
24) 2-Nitrophenol-d4	9.593	143	422364	48.720	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.134	165	762978	35.619	ng/u1	0.00
31) 4-Chloroaniline-d4	10.651	131	1233354	35.274	ng/u1	0.00
46) Dimethylphthalate-d6	13.786	166	2154912	36.285	ng/u1	-0.01
49) Acenaphthylene-d8	14.057	160	2661357	35.112	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.580	143	351253	37.867	ng/u1	0.00
60) Fluorene-d10	15.369	176	1820202	35.715	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	155238	28.656	ng/u1	-0.01
73) Anthracene-d10	17.233	188	2724468	37.450	ng/u1	-0.01
81) Pyrene-d10	19.480	212	3087819	34.666	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.456	264	3323293	38.651	ng/u1	-0.02

Target Compounds Qvalue

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

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