

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011140.D
 Acq On : 27 Jul 2022 21:43
 Operator : CG/JU
 Sample : PB146443BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK443

Quant Time: Jul 28 00:47:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 00:46:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.658	152	320643	20.000	ng/u1	0.00
20) Naphthalene-d8	10.452	136	1343336	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.322	164	697902	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.086	188	1261871	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.204	240	1189998	20.000	ng/u1	# 0.00
88) Perylene-d12	23.539	264	1289071	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	51528	5.793	ng/uL	0.00
4) Pyridine-d5	3.570	84	711050	30.930	ng/u1	0.00
7) Phenol-d5	6.834	99	907006	34.460	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	659152	38.946	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	746242	35.088	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	715365	33.950	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	371883	37.988	ng/u1	0.00
24) 2-Nitrophenol-d4	9.540	143	357251	36.880	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	582975	32.908	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	948812	32.488	ng/u1	0.00
46) Dimethylphthalate-d6	13.740	166	1720165	36.063	ng/u1	0.00
49) Acenaphthylene-d8	14.010	160	2190592	38.924	ng/u1	0.00
54) 4-Nitrophenol-d4	14.540	143	225234	23.331	ng/u1	0.00
60) Fluorene-d10	15.322	176	1427033	35.197	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	166771	26.709	ng/u1	0.00
73) Anthracene-d10	17.186	188	2023661	37.723	ng/u1	0.00
81) Pyrene-d10	19.439	212	2225128	35.603	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.392	264	2316368	37.413	ng/u1	0.00

Target Compounds Qvalue

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

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